

Heat Simulation using MPI

Group 6

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Heat Simulation using MPI

Utilized Functions

Creating a 2d cartesian topology:

```
int MPI_Cart_create(MPI_Comm comm_old, int ndims, int *dims, int *periods,  
                   int reorder, MPI_Comm *comm_cart)
```

Converting between coordinates and rank:

```
int MPI_Cart_rank(MPI_Comm comm, int *coords, int *rank)  
int MPI_Cart_coords(MPI_Comm comm, int rank, int maxdims, int *coords)
```

Nonblocking send and receive commands:

```
int MPI_Isend(void *buf, int count, MPI_Datatype datatype, int dest, int tag,  
             MPI_Comm comm, MPI_Request *request)  
int MPI_Irecv(void *buf, int count, MPI_Datatype datatype, int source,  
             int tag, MPI_Comm comm, MPI_Request *request)  
int MPI_Wait(MPI_Request *request, MPI_Status *status)
```

Send and receive buffers to transfer columns of the matrix

```
// Example: Receiving the left border  
for (i = 1; i < size_y - 1; i++)  
    u[i*size_x] = recvbuf_left[i-1];
```

Heat Simulation using MPI

Code Jacobi

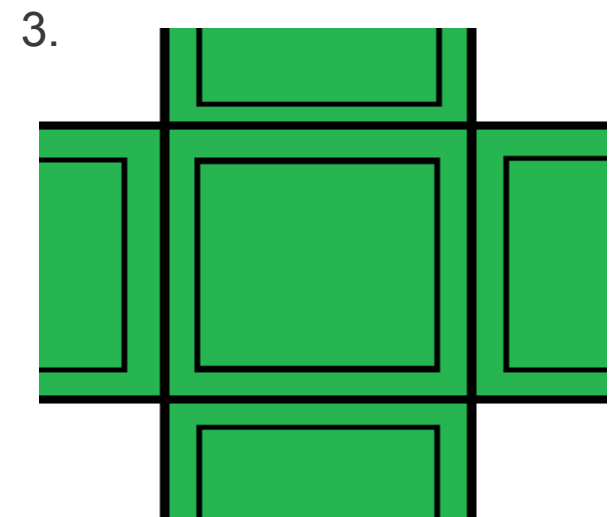
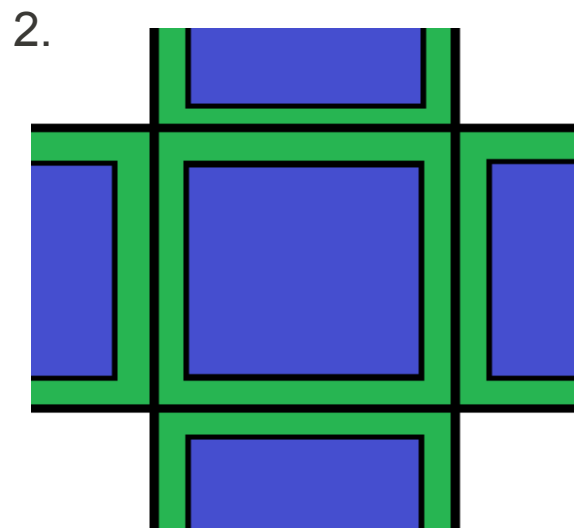
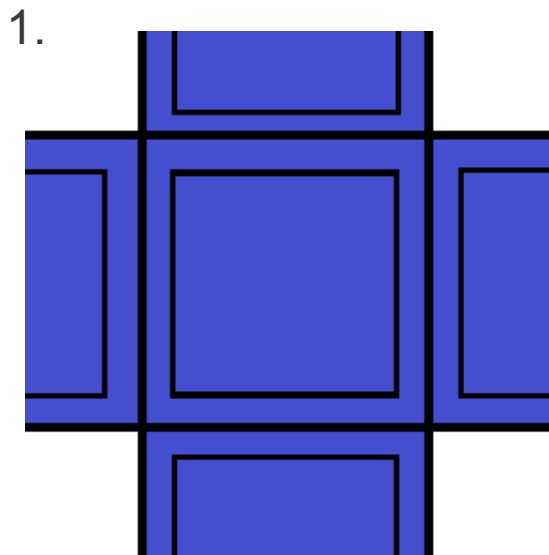
For each iteration do:

- Calculate the borders

- Send border values to neighbors

- Calculate the inner part

- Receive border values from neighbors



Heat Simulation using MPI

Code 1D-Gauss

For each iteration do:

Receive border values from left neighbor

Calculate left border

Send border values to left neighbor

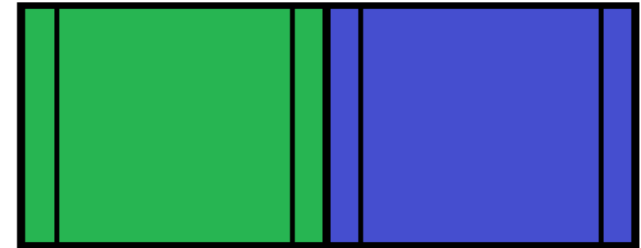
Calculate inner part

Receive border values from right neighbor

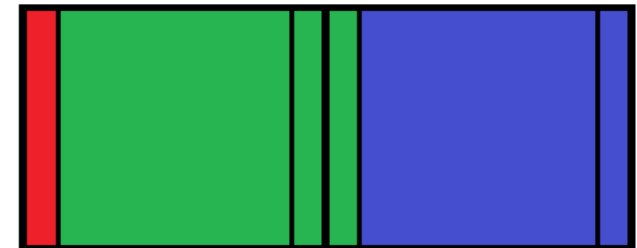
Calculate right border

Send border values to right neighbor

1.



2.



3.



4.



Heat Simulation using MPI

Code 2D-Gauss

For each iteration do:

Receive border values from top and left neighbor

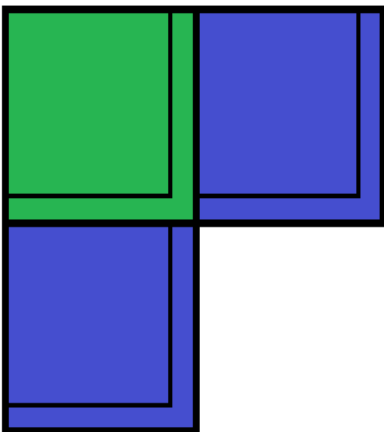
Calculate top and left border and inner part

Receive border values from right and bottom neighbor

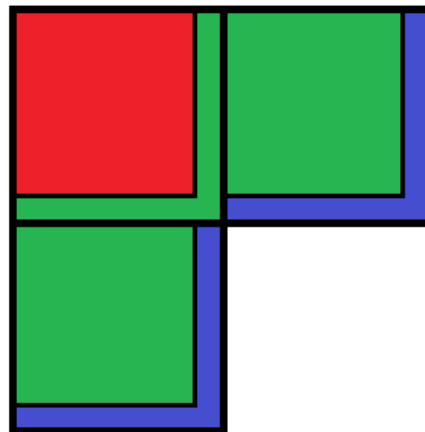
Calculate bottom and right border

Send border values to all neighbors

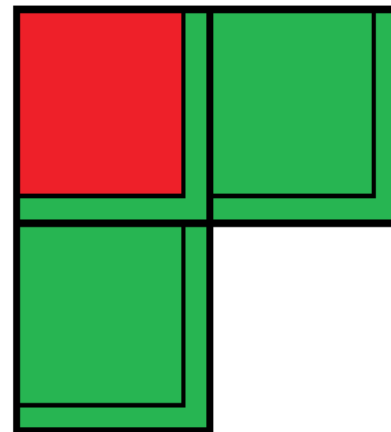
1.



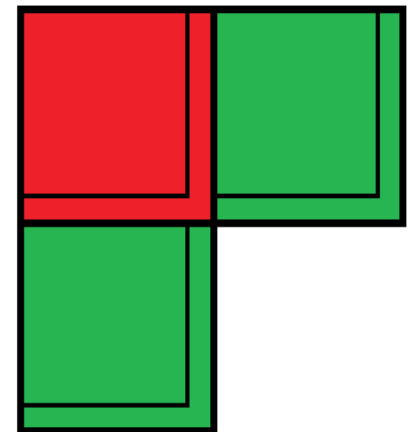
2.



3.



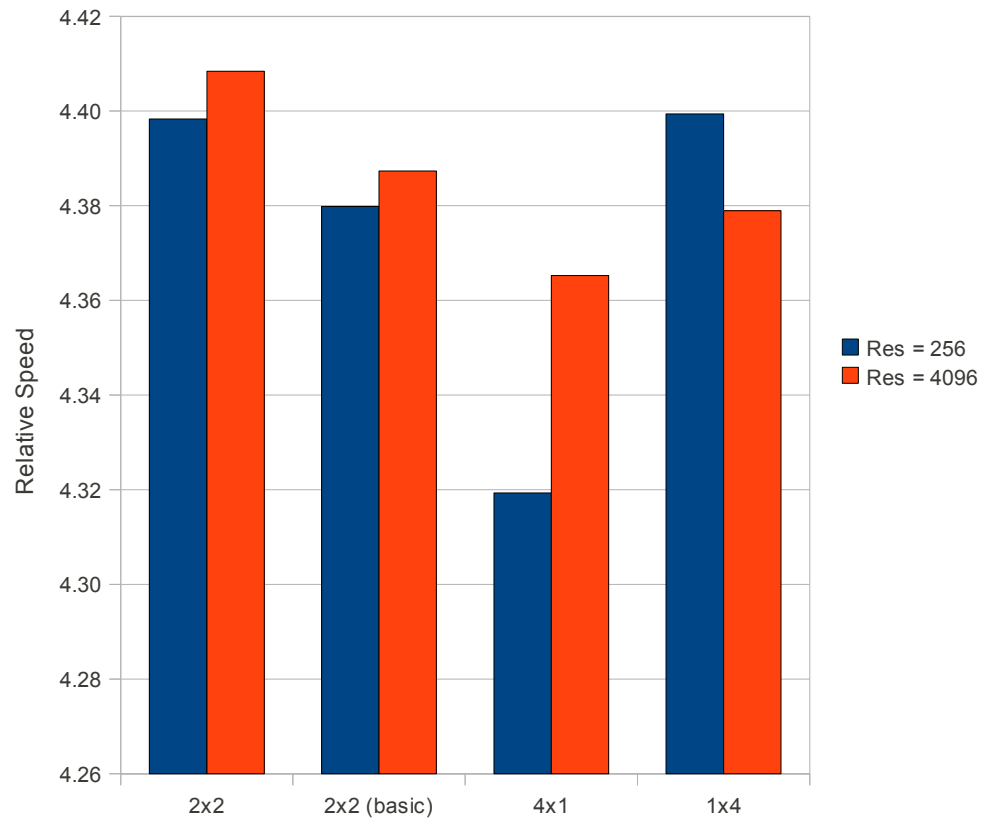
4.



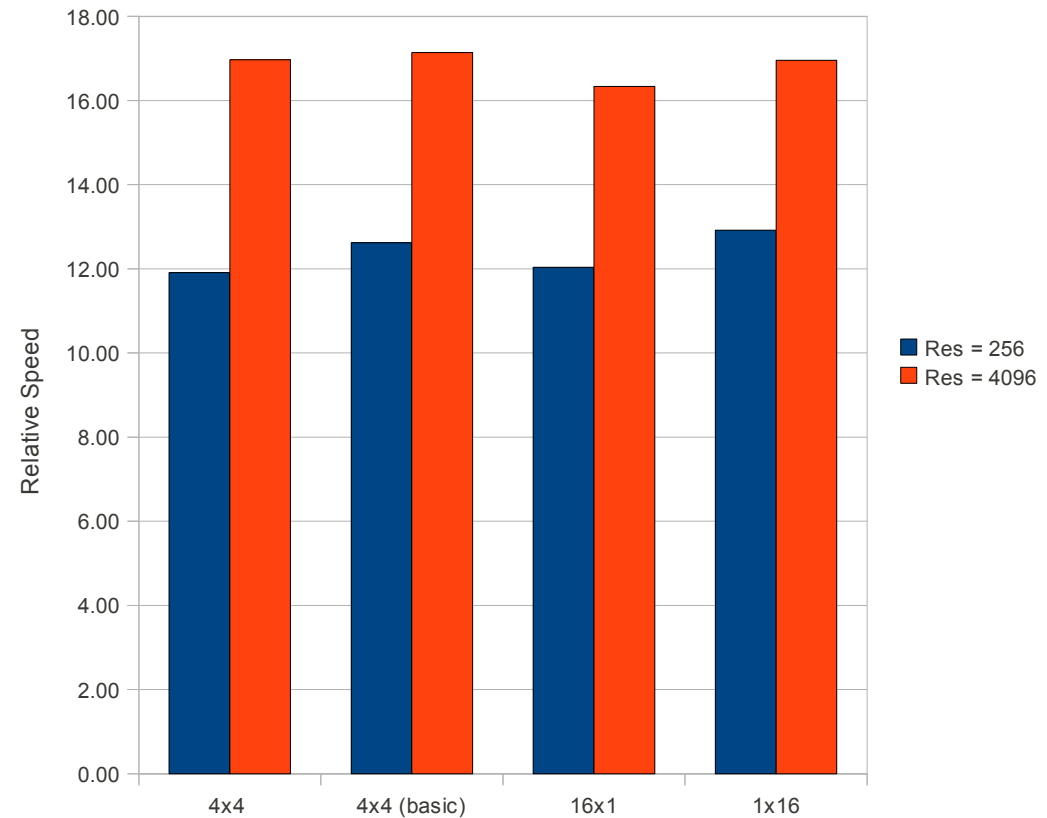
Heat Simulation using MPI

Results Jacobi

4 Threads Jacobi
(lce1)



16 Threads Jacobi
(lce1)

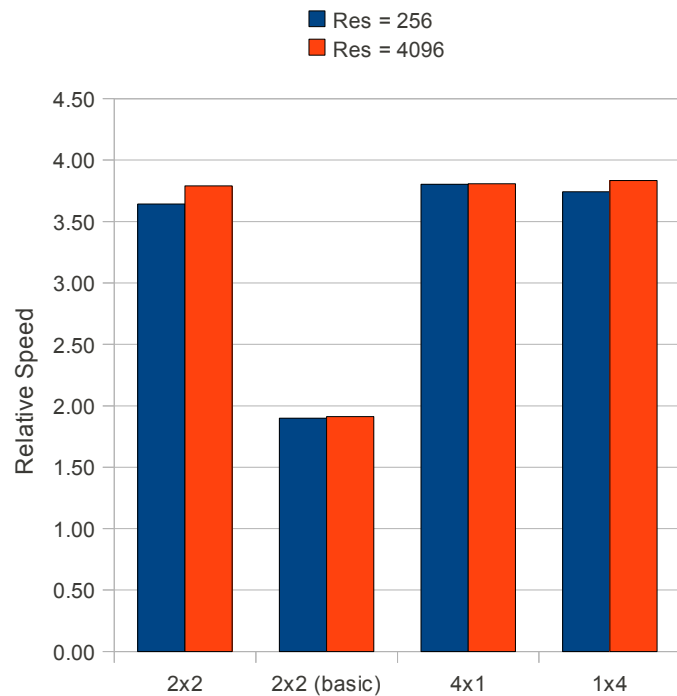


„Basic“ means that the blocks are not divided into inner part and border, but calculated at once.

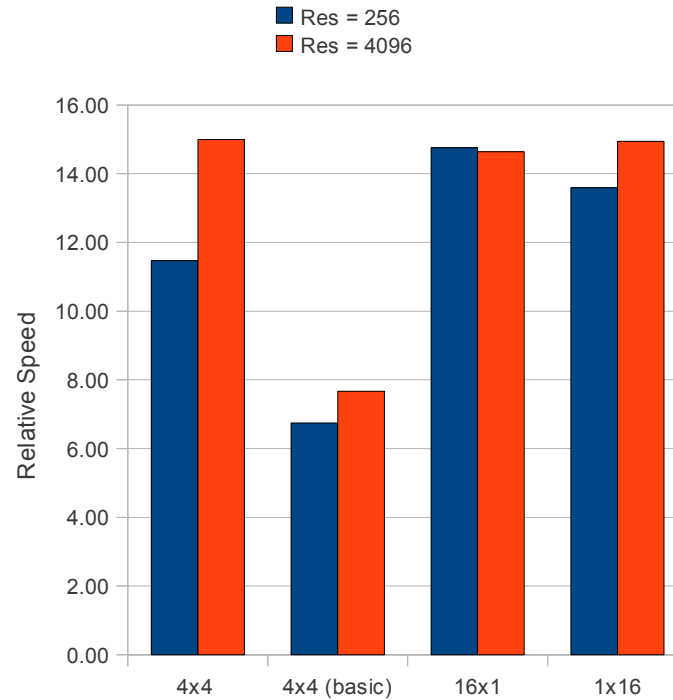
Heat Simulation using MPI

Results Gauss

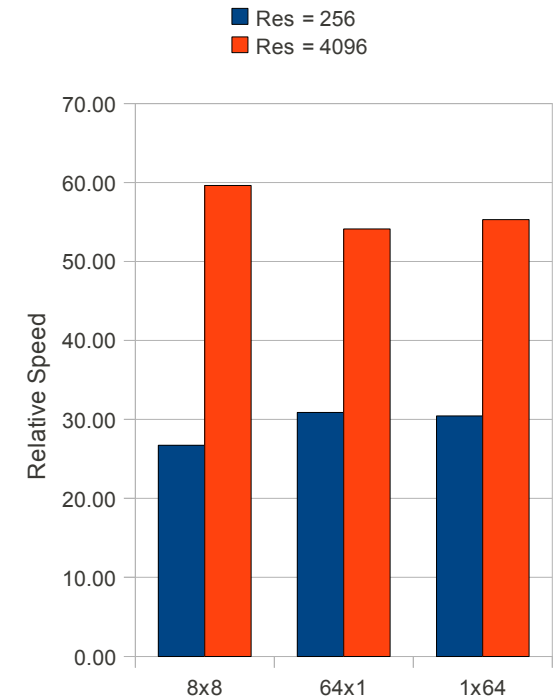
4 Threads Gauss-Seidel
(Ice1)



16 Threads Gauss-Seidel
(Ice1)



64 Threads Gauss-Seidel
(Ice1)



„Basic“ means that the blocks are not divided into upper left part and lower right border, but calculated at once.

Raw Results in MFlops for Gauss (Ice1)

4 Threads

Resolution	sequential	2x2	2x2 (basic)	4x1	1x4
256	842	3068	1600	3202	3152
4096	819	3104	1567	3118	3139

16 Threads

Resolution	sequential	4x4	4x4 (basic)	16x1	1x16
256	842	9657	5680	12425	11443
4096	819	12281	6281	11987	12234

64 Threads

Resolution	sequential	8x8	64x1	1x64
256	842	22521	26017	25639
4096	819	48842	44315	45296

Raw Results in MFlops for Jacobi (Ice1)

4 Threads

Resolution	sequential	2x2	2x2 (basic)	4x1	1x4
256	974	4284	4266	4207	4285
4096	950	4188	4168	4147	4160

16 Threads

Resolution	sequential	4x4	4x4 (basic)	16x1	1x16
256	974	11603	12293	11726	12584
4096	950	16122	16283	15517	16111